

AMENDMENT

Subject matter to be added is in bold and underlined.

Subject matter to be deleted is in bold and strikethrough.

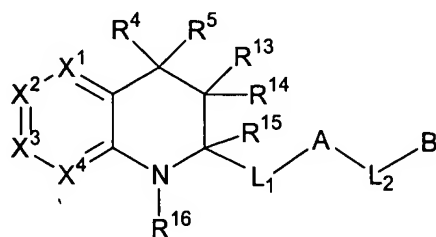
In the Claims:

Please amend Claims 1-3 and enter new claims 24-43 as follows. Please withdraw Claims 11-12 and cancel Claims 15-23 without prejudice to presentation in a Divisional or Continuation application.

This listing of claims will replace all prior versions and listings of claims in the application.

Listing of Claims:

1. (Currently amended) A compound of Formula (I):



(I)

or a stereoisomer or a pharmaceutically acceptable salt or hydrate thereof, wherein:

L_1 is a bond, ~~CH_2~~ , ~~CH_2CH_2~~ , ~~CH_2O~~ , ~~$\text{CH}_2\text{S}(\text{O})_p$~~ , ~~$\text{CH}_2\text{NR}^{10}$~~ , ~~$\text{CH}_2\text{C}(\text{O})$~~ , or ~~$\text{CONR}^{10}$~~ ;

L_2 is a bond, ~~$(\text{CR}^6\text{R}^{6a})_{1-2}$~~ , ~~$\text{CH}_2$~~ , or ~~$\text{O}$~~ , ~~$\text{NR}^{10}$~~ , ~~$\text{C}(\text{O})$~~ , ~~$\text{S}(\text{O})_p$~~ , ~~$(\text{CR}^6\text{R}^{6a})\text{C}(\text{O})$~~ , ~~$\text{C}(\text{O})(\text{CR}^6\text{R}^{6a})$~~ , ~~$(\text{CR}^6\text{R}^{6a})\text{O}$~~ , ~~$\text{O}(\text{CR}^6\text{R}^{6a})$~~ , ~~$(\text{CR}^6\text{R}^{6a})\text{NR}^{10}$~~ , ~~$\text{NR}^{10}(\text{CR}^6\text{R}^{6a})$~~ , ~~$(\text{CR}^6\text{R}^{6a})\text{S}(\text{O})_p$~~ , ~~$\text{S}(\text{O})_p(\text{CR}^6\text{R}^{6a})$~~ , ~~$\text{C}(\text{O})\text{O}$~~ , ~~$\text{OC}(\text{O})$~~ , ~~$\text{C}(\text{O})\text{NR}^8$~~ , ~~$\text{NR}^8\text{C}(\text{O})$~~ , ~~$\text{S}(\text{O})\text{NR}^8$~~ , ~~$\text{S}(\text{O})_2\text{NR}^8$~~ , ~~$\text{NR}^8\text{S}(\text{O})$~~ , or ~~$\text{NR}^8\text{S}(\text{O})_2$~~ ;

A is ~~C_{3-10}~~ **carbocycle** phenyl substituted with 0-3 R^{11} and 0-1 R^{12} , ~~or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$, and substituted 0-3 R^{11} and 0-1 R^{12} ;~~

B is ~~C₁₋₆ alkyl substituted with 0-2 R¹¹ and 0-1 R¹², C₂₋₆ alkenyl substituted with 0-2 R¹¹ and 0-1 R¹², C₂₋₆ alkynyl substituted with 0-2 R¹¹ and 0-1 R¹², C₃₋₁₀ carbocycle~~ phenyl substituted with 0-3 R¹¹ and 0-1 R¹², ~~or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R¹¹ and 0-1 R¹²;~~

X¹, X², X³ and X⁴ independently represent CR¹, CR², or CR³ ~~or N~~;

R¹ is H, -NH₂, -NH(C₁₋₃ alkyl), -N(C₁₋₃ alkyl)₂, -C(=NH)NH₂, -NHC(=NH)NH₂, -C(O)NH₂, -CH₂NH₂, -CH₂NH(C₁₋₃ alkyl), -CH₂N(C₁₋₃ alkyl)₂, -CH₂CH₂NH₂, -CH₂CH₂NH(C₁₋₃ alkyl), -CH₂CH₂N(C₁₋₃ alkyl)₂, -C(=NR⁸)NR⁷R⁹, -NHC(=NR⁸)NR⁷R⁹, -ONHC(=NR⁸)NR⁷R⁹, -NR⁸CH(=NR⁷), -C(=NR^{8a})NR⁷R⁹, -NHC(=NR^{8a})NR⁷R⁹, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -S(O)_pNR⁸R⁹, F, Cl, Br, I, OCF₃, CF₃, OR^a, SR^a, CN or C₁₋₆ alkyl substituted with 1 R^{1a};

R^{1a} is -C(=NR⁸)NR⁷R⁹, -NHC(=NR⁸)NR⁷R⁹, -ONHC(=NR⁸)NR⁷R⁹, -NR⁸CH(=NR⁷), -NR⁷R⁸, -C(O)NR^{7a}R⁸, -S(O)_pNR⁸R⁹, F, OCF₃, CF₃, OR^a, SR^a, or CN;

R² is H, F, Cl, Br, I, OCF₃, CF₃, OR^a, SR^a, CN, NO₂, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹, -S(O)R^c, -S(O)₂R^c, C₁₋₆ alkyl substituted with 0-2 R^{2a}, C₂₋₆ alkenyl substituted with 0-2 R^{2a}, C₂₋₆ alkynyl substituted with 0-2 R^{2a}, or -(CH₂)_r- C₃₋₁₀ carbocycle substituted with 0-3 R^{2b}, ~~or (CH₂)_r- 5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^{2b};~~

each R^{2a} is, independently at each occurrence, H, F, OCF₃, CF₃, OR^a, SR^a, CN, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹, -S(O)R^c, or -S(O)₂R^c;

each R^{2b} is, independently at each occurrence, H, F, Cl, Br, I, OR^a, SR^a, CN, NO₂, CF₃, -SO₂R^c, -NR⁷R⁸, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyloxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(O)-, or C₁₋₄ alkyl-C(O)NH-;

alternately, when R¹ and R² are substituted on adjacent ring carbon atoms, they can be taken together with the ring carbon atoms to which they are attached to form a 5-7 membered carbocycle ~~or heterocycle~~ substituted with 0-2 R^{2b};

R³ is H, F, Cl, Br, I, OCF₃, CF₃, OR^a, SR^a, CN, NO₂, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹, -S(O)R^c, -S(O)₂R^c, C₁₋₆ alkyl substituted with 0-2 R^{3a}, C₂₋₆ alkenyl substituted with 0-2 R^{3a}, C₂₋₆ alkynyl substituted with 0-2 R^{3a}, or -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-3 R^{3b}, ~~or -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^{3b};~~

each R^{3a} is, independently at each occurrence, H, F, OCF₃, CF₃, OR^a, SR^a, CN, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹, -S(O)R^c, or -S(O)₂R^c;

each R^{3b} is, independently at each occurrence, H, F, Cl, Br, I, OR^a, SR^a, CN, NO₂, CF₃, -SO₂R^c, -NR⁷R⁸, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyloxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(O)-, or C₁₋₄ alkyl-C(O)NH-;

R⁴ is ~~H, F, OR^a, SR^a, -NR⁷R⁸, -NR¹⁰C(O)NR^{7a}R⁸, -NR¹⁰SO₂R^c, -C(O)OR^a, -C(O)NR^{7a}R⁸, C₁₋₄ haloalkyl, C₁₋₆ alkyl substituted with 0-3 R^{4a}, C₂₋₆ alkenyl substituted with 0-3 R^{4a}, C₂₋₆ alkynyl substituted with 0-3 R^{4a}, -(CH₂)_r-C₃₋₁₀ carbocycle phenyl substituted with 0-3 R^{4b}, or -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^{4b};~~

~~each R^{4a} is, independently at each occurrence, H, C₁₋₄ alkyl, OR^a, F, =O, CF₃, CN, -C(O)R^a, -C(O)OR^a, -C(O)NR^{7a}R⁸, -NR¹⁰COR^c, or -S(O)_pR^b;~~

each R^{4b} is, independently at each occurrence, H, OH, Cl, F, Br, I, CN, NO₂, CF₃, -C(O)OR^a, -SO₂R^c, -NR⁷R⁸, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyloxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(O)-, C₁₋₄ alkyl-C(O)NH-, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -NR¹⁰S(O)₂NR⁸R⁹, or -S(O)₂NR⁸R⁹;

R^5 is H, F, C_{1-4} haloalkyl, or C_{1-6} alkyl substituted with 0-3 R^{5a} , ~~C_{2-6} alkenyl substituted with 0-3 R^{5a} , C_{2-6} alkynyl substituted with 0-3 R^{5a} , $(CH_2)_r$ C_{3-10} carbocycle substituted with 0-3 R^{5b} , or $(CH_2)_r$ 5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^{5b} ;~~

each R^{5a} is, independently at each occurrence, H, C_{1-4} alkyl, OR^a , F, =O, CF_3 , CN, $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^{7a}R^8$, or $-S(O)_pR^c$;

~~each R^{5b} is, independently at each occurrence, H, OH, Cl, F, Br, I, CN, NO_2 , CF_3 , $-C(O)OR^a$, SO_2R^e , NR^7R^8 , C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-6} cycloalkyl, C_{1-4} haloalkyl, C_{1-4} haloalkyloxy, C_{1-4} alkyloxy, C_{1-4} alkylthio, C_{1-4} alkyl $-C(O)-$, or C_{1-4} alkyl $-C(O)NH-$;~~

each R^6 is, independently at each occurrence, H, C_{1-4} alkyl, $-(CH_2)_rC(O)OR^a$, $-(CH_2)_rS(O)_2NR^{7a}R^8$, or $-(CH_2)_rOR^a$;

each R^{6a} is, independently at each occurrence, H or C_{1-4} alkyl;

each R^7 is, independently at each occurrence, H, C_{1-6} alkyl, $-(CH_2)_n$ -phenyl, $(C_{1-6}$ alkyl) $C(O)-$, $(C_{6-10}$ aryl)- C_{0-4} alkyl- $C(O)-$, $(C_{3-6}$ cycloalkyl)- C_{0-4} alkyl- $C(O)-$, ~~$(5-10$ membered heteroaryl)- C_{0-4} alkyl- $C(O)-$,~~ $(C_{1-4}$ alkyl) $OC(O)-$, $(C_{6-10}$ aryl)- C_{0-4} alkyl- $OC(O)-$, $(C_{1-4}$ alkyl)- $C(O)O-(C_{1-4}$ alkyl)- $OC(O)-$, $(C_{6-10}$ aryl)- $C(O)O-(C_{1-4}$ alkyl)- $OC(O)-$, ~~$(5-10$ membered heteroaryl)- $CH_2-OC(O)-$,~~ $(C_{1-6}$ alkyl)- $NHC(O)-$, $(C_{6-10}$ aryl)- C_{0-4} alkyl- $NHC(O)-$, ~~$(5-10$ membered heteroaryl)- C_{0-4} alkyl- $NHC(O)-$,~~ $(C_{1-6}$ alkyl)- $S(O)_2-$, $(C_{6-10}$ aryl)- $(C_{0-4}$ alkyl)- $S(O)_2-$, ~~$(5-10$ membered heteroaryl)- C_{0-4} alkyl- $S(O)_2-$,~~ $(C_{1-6}$ alkyl) $_2NC(O)-$, phenyl- $NHC(O)-$, benzyl- $NHC(O)-$, $(phenyl)(C_{1-6}$ alkyl) $NC(O)-$, or $(benzyl)(C_{1-6}$ alkyl) $NC(O)-$, wherein said phenyl, and aryl and heteroaryl are substituted with 0-2 R^f ;

each R^{7a} is, independently at each occurrence, H, C_{1-4} alkyl substituted with 0-2 R^{7b} or 0-2 R^{7c} , or $-(CH_2)_r$ C_{3-10} carbocycle substituted with 0-3 R^f , or a $-(CH_2)_r$ 5-12

~~membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted 0-3 R^f;~~

each R^{7b} is, independently at each occurrence, =O, OR^g, F, Cl, Br, I, CN, NO₂, -NR⁷R⁸, -C(O)R^g, -C(O)OR^g, -NR⁸C(O)R^g, -C(O)NR⁸R⁹, -NR⁸C(O)NR⁸R⁹, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl, -S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, or -(CF₂)_rCF₃;

each R^{7c} is, ~~independently at each occurrence~~, C₃₋₁₀ carbocycle substituted with 0-3 R^f; ~~or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted 0-3 R^f;~~

each R⁸ is, independently at each occurrence, H, C₁₋₆ alkyl, or -(CH₂)_n-phenyl;

each R^{8a} is, independently at each occurrence, H, OH, C₁₋₆ alkyl, -(CH₂)_n-phenyl, (C₁₋₆ alkyl)C(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-C(O)-, (C₃₋₆ cycloalkyl)-C₀₋₄ alkyl-C(O)-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-C(O)-~~, (C₁₋₄ alkyl)OC(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-OC(O)-, (C₁₋₄ alkyl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-, (C₆₋₁₀ aryl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-, ~~(5-10 membered heteroaryl)-CH₂-OC(O)-~~, (C₁₋₆ alkyl)-NHC(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-NHC(O)-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-NHC(O)-~~, (C₁₋₆ alkyl)-S(O)₂-, (C₆₋₁₀ aryl)-(C₀₋₄ alkyl)-S(O)₂-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-S(O)₂-~~, C₁₋₄ alkoxy, (C₁₋₄ alkyl)C(O)O-, (C₆₋₁₀ aryl)-(C₀₋₄ alkyl)-C(O)O-; wherein said phenyl, and aryl ~~and heteroaryl~~ are substituted with 0-2 R^f;

~~alternatively, R⁷ and R⁸, or R^{7a} and R⁸, when attached to the same nitrogen, combine to form a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;~~

each R⁹ is, independently at each occurrence, H, C₁₋₆ alkyl, or -(CH₂)_n-phenyl;

each R^{10} is, independently at each occurrence, H, C_{1-6} alkyl substituted with 0-2 R^{10a} , C_{2-6} alkenyl substituted with 0-2 R^{10a} , C_{2-6} alkynyl substituted with 0-2 R^{10a} , $(C_{1-6} \text{ alkyl})C(O)-$, $(C_{3-6} \text{ cycloalkyl})C_{1-3} \text{ alkyl}-C(O)-$, $(C_{3-6} \text{ cycloalkyl})C(O)-$, phenyl- $C(O)-$, benzyl- $C(O)-$, benzyl- $S(O)_2-$, $(C_{1-6} \text{ alkyl})NHC(O)-$, $(C_{1-6} \text{ alkyl})_2NC(O)-$, phenyl- $NHC(O)-$, benzyl- $NHC(O)-$, (phenyl) $(C_{1-6} \text{ alkyl})NC(O)-$, (benzyl) $(C_{1-6} \text{ alkyl})NC(O)-$, $(C_{1-6} \text{ alkyl})-S(O)_2-$, phenyl- $S(O)_2-$, or $-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-3 R^d , ~~or $-(CH_2)_r$ 5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^d ;~~

each R^{10a} is, independently at each occurrence, H, C_{1-4} alkyl, OR^a , Cl, F, Br, I, $=O$, CF_3 , CN, NO_2 , $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^7aR^8$, or $-S(O)_pR^c$;

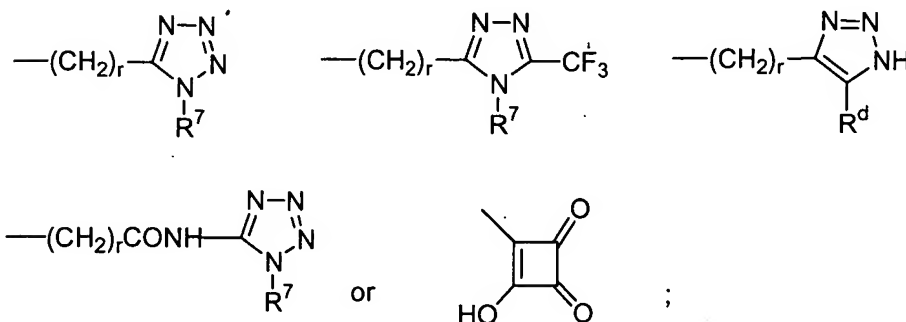
each R^{11} is, independently at each occurrence, H, $=O$, $-(CH_2)_r-OR^a$, F, Cl, Br, I, CN, NO_2 , $-(CH_2)_r-NR^7R^8$, $-C(O)R^a$, $-C(O)OR^a$, $-NR^8C(O)R^a$, $-NR^8C(O)OR^a$, $-C(O)NR^7aR^8$, $-NR^8C(O)NR^8R^9$, $-SO_2NR^8R^9$, $-NR^8SO_2NR^8R^9$, $-NR^8SO_2-C_{1-4} \text{ alkyl}$, $-NR^8SO_2CF_3$, $-NR^8SO_2\text{-phenyl}$, $-S(O)_2CF_3$, $-S(O)_p-C_{1-4} \text{ alkyl}$, $-S(O)_p\text{-phenyl}$, $-(CF_2)_rCF_3$, $C_{1-6} \text{ alkyl}$ substituted with 0-2 R^{11a} , $C_{2-6} \text{ alkenyl}$ substituted with 0-2 R^{11a} , $C_{2-6} \text{ alkynyl}$ substituted with 0-2 R^{11a} , $C_{1-6} \text{ alkyl}$ substituted with 0-2 R^{11b} , $C_{2-6} \text{ alkenyl}$ substituted with 0-2 R^{11b} , or $C_{2-6} \text{ alkynyl}$ substituted with 0-2 R^{11b} ;

each R^{11a} is, independently at each occurrence, $=O$, OR^a , F, Cl, Br, I, CN, NO_2 , $-NR^7R^8$, $-C(O)R^a$, $-C(O)OR^a$, $-NR^8C(O)R^a$, $-C(O)NR^7aR^8$, $-NR^8C(O)NR^8R^9$, $-SO_2NR^8R^9$, $-NR^8SO_2NR^8R^9$, $-NR^8SO_2-C_{1-4} \text{ alkyl}$, $-NR^8SO_2CF_3$, $-NR^8SO_2\text{-phenyl}$, $-S(O)_2CF_3$, $-S(O)_p-C_{1-4} \text{ alkyl}$, $-S(O)_p\text{-phenyl}$, or $-(CF_2)_rCF_3$;

each R^{11b} is, independently at each occurrence, C_{3-10} carbocycle substituted with 0-3 R^d ; or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted 0-3 R^d ;

each R^{12} is, independently at each occurrence, OR^{12a} , $-C(O)NR^7aR^8$, $-(CH_2)_rCO_2R^{12a}$, $-(CH_2)_rSO_3H$, $-OSO_3H$, $-(CH_2)_rPO_3H$, $-OPO_3H_2$, $-PO_3H_2$,

-NHPO₃H₂, -NHCOCF₃, -NHSO₂CF₃, -CONHNHSO₂CF₃, -C(CF₃)₂OH, -SO₂NHR^{12a},
-CONHSO₂NHR^{12a}, -SO₂NHCOR^{12a}, -SO₂NHCO₂R^{12a}, -CONHSO₂R^{12b},
-NHSO₂R^{12b}, -CONHOR^{12b},



each R^{12a} is, independently at each occurrence, H, C₁₋₆ alkyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-3 R^d; or -(CH₂)_r-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^d;

each R^{12b} is, independently at each occurrence, C₁₋₆ alkyl substituted with 0-2 R^{12c}, C₂₋₆ alkenyl substituted with 0-2 R^{12c}, C₂₋₆ alkynyl substituted with R^{12c}, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-3 R^{12c}, or -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^{12c};

each R^{12c} is, independently at each occurrence, H, F, Cl, Br, I, CF₃, OCF₃, CN, NO₂, OR^a, -CO₂R^a, -NR⁷R⁸, -SO₂R^c, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-3 R^d; or -(CH₂)_r-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^d;

R¹³ is H, C₁₋₄ alkyl, (NR⁷R⁸)C₁₋₄ alkyl, (SR^c)C₁₋₄ alkyl, (OR^a)C₁₋₄ alkyl, OR^a, F, CF₃, -C(O)R^a, -C(O)OR^a, -C(O)NR^{7a}R⁸, or -S(O)_pR^c;

R¹⁴ is H, C₁₋₄ alkyl, (NR⁷R⁸)C₁₋₄ alkyl, (SR^c)C₁₋₄ alkyl, (OR^a)C₁₋₄ alkyl, OR^a, F, CF₃, -C(O)R^a, -C(O)OR^a, -C(O)NR^{7a}R⁸, or -S(O)_pR^c;

alternately, R¹³ and R¹⁴ may be taken together to be =O;

R¹⁵ is H or C₁₋₄ alkyl;

R¹⁶ is H, C₁₋₄ alkyl, benzyl, C₁₋₄ alkyl-C(O)-, C₁₋₄ alkyl-S(O)₂-, or C₁₋₄ alkyl-OC(O)-;

each R^a is, independently at each occurrence, H, C₁₋₄ alkyl, -(CH₂)_r-CO₂R^g, -(CH₂)_r-C₃₋₇ cycloalkyl, or -(CH₂)_r-C₆₋₁₀ aryl, ~~or -(CH₂)_r-5-10 membered heteroaryl,~~ wherein said aryl ~~or heteroaryl groups are~~ is substituted with 0-2 R^f;

each R^b is, independently at each occurrence, CF₃, OH, C₁₋₄ alkoxy, C₁₋₆ alkyl, or -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^d, ~~or -(CH₂)_r-5-10 membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p and substituted with 0-2 R^d;~~

each R^c is, independently at each occurrence, C₁₋₄ alkyl, C₆₋₁₀ aryl, or ~~5-10 membered heteroaryl,~~ (C₆₋₁₀ aryl)-C₁₋₄ alkyl, ~~or (5-10 membered heteroaryl)-C₁₋₄ alkyl,~~ wherein said aryl ~~and heteroaryl groups are~~ is substituted with 0-2 R^d;

each R^d is, independently at each occurrence, H, =O, OR^a, F, Cl, Br, I, CN, NO₂, -NR⁷R⁸, -C(O)R^a, -C(O)OR^a, -NR⁸C(O)R^a, -C(O)NR^{7a}R⁸, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl, -S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, -(CF₂)_rCF₃, C₁₋₆ alkyl substituted with 0-2 R^e, C₂₋₆ alkenyl substituted with 0-2 R^e, or C₂₋₆ alkynyl substituted with 0-2 R^e;

each R^e is, independently at each occurrence, =O, OR^a, F, Cl, Br, I, CN, NO₂, -NR⁸R⁹, -C(O)R^a, -C(O)OR^a, -NR⁸C(O)R^a, -C(O)NR^{7a}R⁸, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl, -S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, or -(CF₂)_rCF₃;

each R^f is, independently at each occurrence, H, =O, OR^g, F, Cl, Br, I, CN, NO₂, -NR⁸R⁹, -C(O)R^g, -C(O)OR^g, -NR⁸C(O)R^g, -C(O)NR⁸R⁹, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl, -S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, -(CF₂)_rCF₃, C₁₋₆ alkyl, C₂₋₆ alkenyl, or C₂₋₆ alkynyl;

each R^g is, independently at each occurrence, H, C₁₋₆ alkyl, or

-(CH₂)_n-phenyl;

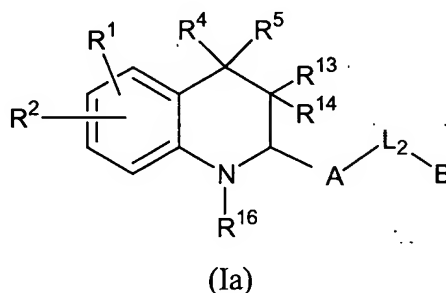
n, at each occurrence, is selected from 0, 1, 2, 3, and 4;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, and 4; and

provided that when L₁ is a bond and A is phenyl or a 6-membered aromatic N-heterocycle, then ring A is not substituted ortho to L₁ with OH, halogen, -CO₂H, -C(O)O-C₁₋₄ alkyl, -O-phenyl, -O-benzyl, -NR⁷R⁸, -CH₂OR^a, haloalkyl, -S-C₁₋₄ alkyl, or -NHSO₂-C₁₋₄ alkyl.

2. (Currently amended) A compound according to Claim 1, wherein the compound is of Formula (Ia):



or a stereoisomer or a pharmaceutically acceptable salt or hydrate thereof, wherein:

~~L₂ is a bond, (CR⁶R^{6a})₁₋₂, O, NR¹⁰, C(O), S(O)_p, CR⁶R^{6a}C(O), C(O)(CR⁶R^{6a}), (CR⁶R^{6a})O, O(CR⁶R^{6a}), (CR⁶R^{6a})NR¹⁰, NR¹⁰(CR⁶R^{6a}), (CR⁶R^{6a})S(O)_p, S(O)_p(CR⁶R^{6a}), C(O)O, OC(O), C(O)NR⁸, NR⁸C(O), S(O)NR⁸, S(O)₂NR⁸, NR⁸S(O), or NR⁸S(O)₂;~~

~~A is phenyl substituted with 0-2 R¹¹ and 0-1 R¹², or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted 0-2 R¹¹ and 0-1 R¹²;~~

~~B is phenyl substituted with 0-2 R¹¹ and 0-1 R¹², or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R¹¹ and 0-1 R¹²;~~

R¹ is H, -NH₂, -NH(C₁₋₃ alkyl), -N(C₁₋₃ alkyl)₂, -C(=NH)NH₂,

-NHC(=NH)NH₂, -C(O)NH₂, -CH₂NH₂, -CH₂NH(C₁-C₃ alkyl),
-CH₂N(C₁-C₃ alkyl)₂, -CH₂CH₂NH₂, -CH₂CH₂NH(C₁-C₃ alkyl),
-CH₂CH₂N(C₁-C₃ alkyl)₂, -C(=NR⁸)NR⁷R⁹, -NHC(=NR⁸)NR⁷R⁹,
-ONHC(=NR⁸)NR⁷R⁹, -NR⁸CH(=NR⁷), -C(=NR^{8a})NR⁷R⁹, -NHC(=NR^{8a})NR⁷R⁹,
-NR⁷R⁸, -C(O)NR^{7a}R⁸, -S(O)_pNR⁸R⁹, F, Cl, Br, I, OCF₃, CF₃, OR^a, SR^a, CN or
C₁₋₆ alkyl substituted with 1 R^{1a};

R^{1a} is -C(=NR⁸)NR⁷R⁹, -NHC(=NR⁸)NR⁷R⁹, -ONHC(=NR⁸)NR⁷R⁹,
-NR⁸CH(=NR⁷), -NR⁷R⁸, -C(O)NR^{7a}R⁸, -S(O)_pNR⁸R⁹, F, OCF₃, CF₃, OR^a, SR^a, or
CN;

R² is H, F, OR^a, CN, -NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹,
-S(O)R^c, -S(O)₂R^c, C₁₋₆ alkyl substituted with 0-2 R^{2a}, or -(CH₂)_r-C₃-C₇ carbocycle
substituted with 0-2 R^{2b}, ~~or -(CH₂)_r-5-7 membered heterocycle consisting of: carbon
atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p,
and substituted with 0-2 R^{2b};~~

each R^{2a} is, independently at each occurrence, H, F, OCF₃, CF₃, OR^a, SR^a, CN,
-NR⁷R⁸, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -S(O)_pNR⁸R⁹, -S(O)R^c, or -S(O)₂R^c;

each R^{2b} is, independently at each occurrence, H, F, OR^a, SR^a, CN, NO₂, CF₃,
-SO₂R^c, -NR⁷R⁸, C₁-C₆ alkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₃-C₆ cycloalkyl, C₁-C₄
haloalkyl, C₁-C₄ haloalkyloxy-, C₁-C₄ alkyloxy-, C₁-C₄ alkylthio-, C₁-C₄ alkyl-C(O)-,
or C₁-C₄ alkyl-C(O)NH-;

alternately, when R¹ and R² are substituted on adjacent ring carbon atoms, they
can be taken together with the ring carbon atoms to which they are attached to form a 5-7
membered carbocycle ~~or heterocycle~~ substituted with 0-2 R^{2b};

R⁴ is ~~H, C₁-C₄ haloalkyl, C₁-C₆ alkyl substituted with 0-3 R^{4a}, C₂-C₆ alkenyl
substituted with 0-3 R^{4a}, C₂-C₆ alkynyl substituted with 0-3 R^{4a}, (CH₂)_r-C₃-C₈
carbocycle phenyl~~ substituted with 0-3 R^{4b}, ~~or -(CH₂)_r-5-6 membered heterocycle
consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting
of N, O, and S(O)_p, and substituted with 0-3 R^{4b};~~

~~each R^{4a} is, independently at each occurrence, H, C₁₋₄ alkyl, OR^a, F, =O, CF₃, CN, C(O)R^a, C(O)OR^a, C(O)NR^{7a}R⁸, NR¹⁰COR^e, or S(O)_pR^b;~~

each R^{4b} is, independently at each occurrence, H, OH, Cl, F, Br, CN, NO₂, CF₃, -C(O)OR^a, -SO₂R^c, -NR⁷R⁸, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyloxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(O)-, C₁₋₄ alkyl-C(O)NH-, -C(O)NR^{7a}R⁸, -NR¹⁰C(O)R^b, -NR¹⁰S(O)₂NR⁸R⁹, or -S(O)₂NR⁸R⁹;

R⁵ is H, F, C₁₋₄ haloalkyl, or C₁₋₆ alkyl substituted with 0-2 R^{5a}, ~~C₂₋₆ alkenyl substituted with 0-2 R^{5a}, C₂₋₆ alkynyl substituted with 0-2 R^{5a}, (CH₂)_r-C₃₋₇ cycloalkyl substituted with 0-2 R^{5b}, (CH₂)_r-phenyl substituted with 0-2 R^{5b}, or (CH₂)_r-5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{5b};~~

~~R^{5a} is, independently at each occurrence, H, C₁₋₄ alkyl, OR^a, F, =O, CF₃, CN, C(O)R^a, C(O)OR^a, C(O)NR^{7a}R⁸, or S(O)_pR^e;~~

~~each R^{5b} is, independently at each occurrence, H, OH, Cl, F, Br, CN, NO₂, CF₃, C(O)OR^a, SO₂R^c, NR⁷R⁸, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₆ cycloalkyl, C₁₋₄ haloalkyl, C₁₋₄ haloalkyloxy-, C₁₋₄ alkyloxy-, C₁₋₄ alkylthio-, C₁₋₄ alkyl-C(O)-, or C₁₋₄ alkyl-C(O)NH-;~~

each R⁶ is, independently at each occurrence, H, C₁₋₄ alkyl, -(CH₂)_rC(O)OR^a, -(CH₂)_rS(O)₂NR^{7a}R⁸, or -(CH₂)_rOR^a;

each R^{6a} is, independently at each occurrence, H or C₁₋₄ alkyl;

each R⁷ is, independently at each occurrence, H, C₁₋₆ alkyl, -(CH₂)_n-phenyl, (C₁₋₆ alkyl)C(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-C(O)-, (C₃₋₆ cycloalkyl)-C₀₋₄ alkyl-C(O)-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-C(O)-~~, (C₁₋₄ alkyl)OC(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-OC(O)-, (C₁₋₄ alkyl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-, (C₆₋₁₀ aryl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-, ~~(5-10 membered heteroaryl)-CH₂-OC(O)-~~, (C₁₋₆ alkyl)-NHC(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-NHC(O)-,

~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-NHC(O)-~~, (C₁₋₆ alkyl)-S(O)₂-,
(C₆₋₁₀ aryl)-(C₀₋₄ alkyl)-S(O)₂-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-S(O)₂-~~,
(C₁₋₆ alkyl)₂NC(O)-, phenyl-NHC(O)-, benzyl-NHC(O)-, (phenyl)(C₁₋₆ alkyl)NC(O)-, or
(benzyl)(C₁₋₆ alkyl)NC(O)-, wherein said phenyl, and aryl ~~and heteroaryl~~ are
substituted with 0-2 R^f;

each R^{7a} is, independently at each occurrence, H, C₁₋₄ alkyl substituted with 0-1
R^{7b} or 0-1 R^{7c}, -(CH₂)_r-C₃₋₇ cycloalkyl substituted with 0-2 R^f, or -(CH₂)_r-phenyl
substituted with 0-3 R^f, ~~or a (CH₂)_r-5-6 membered heterocycle consisting of: carbon
atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p,
and substituted 0-3 R^f;~~

each R^{7b} is, independently at each occurrence, =O, OR^g, F, Cl, Br, I, CN, NO₂,
-NR⁷R⁸, -C(O)R^g, -C(O)OR^g, -NR⁸C(O)R^g, -C(O)NR⁸R⁹, -NR⁸C(O)NR⁸R⁹,
-SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl,
-S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, or -(CF₂)_rCF₃;

each R^{7c} is, ~~independently at each occurrence~~, C₃₋₁₀ carbocycle substituted
with 0-3 R^f; ~~or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4
heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted
0-3 R^f;~~

each R⁸ is, independently at each occurrence, H, C₁₋₆ alkyl, or -(CH₂)_n-phenyl;

each R^{8a} is, independently at each occurrence, H, OH, C₁₋₆ alkyl,
-(CH₂)_n-phenyl, (C₁₋₆ alkyl)C(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-C(O)-,
(C₃₋₆ cycloalkyl)-C₀₋₄ alkyl-C(O)-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-C(O)-~~,
(C₁₋₄ alkyl)OC(O)-, (C₆₋₁₀ aryl)-C₀₋₄ alkyl-OC(O)-,
(C₁₋₄ alkyl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-, (C₆₋₁₀ aryl)-C(O)O-(C₁₋₄ alkyl)-OC(O)-,
~~(5-10 membered heteroaryl)-CH₂-OC(O)-~~, (C₁₋₆ alkyl)-NHC(O)-,
(C₆₋₁₀ aryl)-C₀₋₄ alkyl-NHC(O)-, ~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-NHC(O)-~~,
(C₁₋₆ alkyl)-S(O)₂-, (C₆₋₁₀ aryl)-(C₀₋₄ alkyl)-S(O)₂-,

~~(5-10 membered heteroaryl)-C₀₋₄ alkyl-S(O)₂-~~, C₁₋₄ alkoxy, (C₁₋₄ alkyl)C(O)O-, or (C₆₋₁₀ aryl)-(C₀₋₄ alkyl)-C(O)O-; wherein said phenyl, and aryl and heteroaryl are substituted with 0-2 R^f;

~~alternatively, R⁷ and R⁸, or R^{7a} and R⁸, when attached to the same nitrogen, combine to form a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;~~

each R⁹ is, independently at each occurrence, H, C₁₋₆ alkyl, or -(CH₂)_n-phenyl;

each R¹⁰ is, independently at each occurrence, H, C₁₋₆ alkyl substituted with 0-2 R^{10a}, C₂₋₆ alkenyl substituted with 0-2 R^{10a}, C₂₋₆ alkynyl substituted with 0-2 R^{10a}, (C₁₋₆ alkyl)C(O)-, (C₃₋₆ cycloalkyl)C₁₋₃ alkyl-C(O)-, (C₃₋₆ cycloalkyl)C(O)-, phenyl-C(O)-, benzyl-C(O)-, benzyl-S(O)₂-, (C₁₋₆ alkyl)NHC(O)-, (C₁₋₆ alkyl)₂NC(O)-, phenyl-NHC(O)-, benzyl-NHC(O)-, (phenyl)(C₁₋₆ alkyl)NC(O)-, (benzyl)(C₁₋₆ alkyl)NC(O)-, (C₁₋₆ alkyl)-S(O)₂-, phenyl-S(O)₂-, or -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-3 R^d, ~~or -(CH₂)_r-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-3 R^d;~~

each R^{10a} is, independently at each occurrence, H, C₁₋₄ alkyl, OR^a, Cl, F, Br, I, =O, CF₃, CN, NO₂, -C(O)R^a, -C(O)OR^a, -C(O)NR^{7a}R⁸, or -S(O)_pR^c;

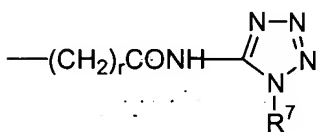
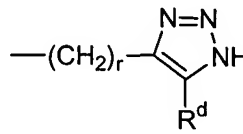
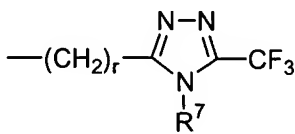
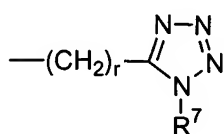
each R¹¹ is, independently at each occurrence, H, =O, -(CH₂)_r-OR^a, F, Cl, Br, I, CN, NO₂, -(CH₂)_r-NR⁷R⁸, -C(O)R^a, -C(O)OR^a, -NR⁸C(O)R^a, -NR⁸C(O)OR^a, -C(O)NR^{7a}R⁸, -NR⁸C(O)NR⁸R⁹, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl, -S(O)₂CF₃, -S(O)_p-C₁₋₄ alkyl, -S(O)_p-phenyl, -(CF₂)_rCF₃; C₁₋₆ alkyl substituted with 0-2 R^{11a}, C₂₋₆ alkenyl substituted with 0-2 R^{11a}, C₂₋₆ alkynyl substituted with 0-2 R^{11a}, C₁₋₆ alkyl substituted with 0-2 R^{11b}, C₂₋₆ alkenyl substituted with 0-2 R^{11b}, or C₂₋₆ alkynyl substituted with 0-2 R^{11b};

each R^{11a} is, independently at each occurrence, =O, OR^a, F, Cl, Br, I, CN, NO₂, -NR⁷R⁸, -C(O)R^a, -C(O)OR^a, -NR⁸C(O)R^a, -C(O)NR^{7a}R⁸, -NR⁸C(O)NR⁸R⁹, -SO₂NR⁸R⁹, -NR⁸SO₂NR⁸R⁹, -NR⁸SO₂-C₁₋₄ alkyl, -NR⁸SO₂CF₃, -NR⁸SO₂-phenyl,

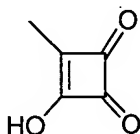
$-\text{S}(\text{O})_2\text{CF}_3$, $-\text{S}(\text{O})_p\text{-C}_{1-4}$ alkyl, $-\text{S}(\text{O})_p\text{-phenyl}$, or $-(\text{CF}_2)_r\text{CF}_3$;

each R^{11b} is, independently at each occurrence, C_{3-10} carbocycle substituted with 0-3 R^d ; or a 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$, and substituted 0-3 R^d ;

each R^{12} is, independently at each occurrence, OR^{12a} , $-\text{C}(\text{O})\text{NR}^7\text{R}^8$, $-(\text{CH}_2)_r\text{CO}_2\text{R}^{12a}$, $-(\text{CH}_2)_r\text{SO}_3\text{H}$, $-\text{OSO}_3\text{H}$, $-(\text{CH}_2)_r\text{PO}_3\text{H}$, $-\text{OPO}_3\text{H}_2$, $-\text{PO}_3\text{H}_2$, $-\text{NHPO}_3\text{H}_2$, $-\text{NHCOCF}_3$, $-\text{NHSO}_2\text{CF}_3$, $-\text{CONHNHSO}_2\text{CF}_3$, $-\text{C}(\text{CF}_3)_2\text{OH}$, $-\text{SO}_2\text{NHR}^{12a}$, $-\text{CONHSO}_2\text{NHR}^{12a}$, $-\text{SO}_2\text{NHCOR}^{12a}$, $-\text{SO}_2\text{NHCO}_2\text{R}^{12a}$, $-\text{CONHSO}_2\text{R}^{12b}$, $-\text{NHSO}_2\text{R}^{12b}$, $-\text{CONHOR}^{12b}$,



or



each R^{12a} is, independently at each occurrence, H, C_{1-6} alkyl, $-(\text{CH}_2)_r\text{-C}_{3-10}$ carbocycle substituted with 0-3 R^d ; or $-(\text{CH}_2)_r\text{-5-10}$ membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$, and substituted with 0-3 R^d ;

each R^{12b} is, independently at each occurrence, $\text{C}_1\text{-C}_6$ alkyl substituted with 0-2 R^{12c} , $\text{C}_2\text{-C}_6$ alkenyl substituted with 0-2 R^{12c} , $\text{C}_2\text{-C}_6$ alkynyl substituted with 0-2 R^{12c} , $-(\text{CH}_2)_r\text{-C}_{3-10}$ carbocycle substituted with 0-3 R^{12c} , or $-(\text{CH}_2)_r\text{-5-10}$ membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $\text{S}(\text{O})_p$, and substituted with 0-3 R^{12c} ;

each R^{12c} is, independently at each occurrence, H, F, Cl, Br, I, CF_3 , OCF_3 , CN, NO_2 , OR^a , $-\text{CO}_2\text{R}^a$, $-\text{NR}^7\text{R}^8$, $-\text{SO}_2\text{R}^c$, $\text{C}_1\text{-C}_6$ alkyl, $\text{C}_2\text{-C}_6$ alkenyl, $\text{C}_2\text{-C}_6$ alkynyl,

$-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-3 R^d ; or $-(CH_2)_r-5-10$ membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^d ;

R^{13} is H or C_{1-4} alkyl;

R^{14} is H or C_{1-4} alkyl;

R^{16} is H, C_{1-4} alkyl, benzyl, C_{1-4} alkyl-C(O)-, C_{1-4} alkyl-S(O)₂-, or C_{1-4} alkyl-OC(O)-;

each R^a is, independently at each occurrence, H, C_{1-4} alkyl, $-(CH_2)_r-CO_2R^g$, $-(CH_2)_r-C_{3-7}$ cycloalkyl, or $-(CH_2)_r-C_{6-10}$ aryl, ~~or $-(CH_2)_r-5-10$ membered heteroaryl,~~ wherein said aryl ~~or heteroaryl groups are~~ is substituted with 0-2 R^f ;

each R^b is, independently at each occurrence, CF_3 , OH, C_{1-4} alkoxy, C_{1-6} alkyl, or $-(CH_2)_r-C_{3-10}$ carbocycle substituted with 0-2 R^d , ~~or $-(CH_2)_r-5-10$ membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^d ;~~

each R^c is, independently at each occurrence, C_{1-4} alkyl, C_{6-10} aryl, ~~5-10 membered heteroaryl, or $(C_{6-10}$ aryl)- C_{1-4} alkyl, or (5-10 membered heteroaryl)- C_{1-4} alkyl, wherein said aryl and heteroaryl groups are substituted with 0-2 R^d ;~~

each R^d is, independently at each occurrence, H, =O, OR^a , F, Cl, Br, I, CN, NO_2 , $-NR^7R^8$, $-C(O)R^a$, $-C(O)OR^a$, $-NR^8C(O)R^a$, $-C(O)NR^7aR^8$, $-SO_2NR^8R^9$, $-NR^8SO_2NR^8R^9$, $-NR^8SO_2-C_{1-4}$ alkyl, $-NR^8SO_2CF_3$, $-NR^8SO_2$ -phenyl, $-S(O)_2CF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$ -phenyl, $-(CF_2)_rCF_3$, C_1-C_6 alkyl substituted with 0-2 R^e , C_2-C_6 alkenyl substituted with 0-2 R^e , or C_2-C_6 alkynyl substituted with 0-2 R^e ;

each R^e is, independently at each occurrence, =O, OR^a , F, Cl, Br, I, CN, NO_2 , $-NR^8R^9$, $-C(O)R^a$, $-C(O)OR^a$, $-NR^8C(O)R^a$, $-C(O)NR^7aR^8$, $-SO_2NR^8R^9$, $-NR^8SO_2NR^8R^9$, $-NR^8SO_2-C_{1-4}$ alkyl, $-NR^8SO_2CF_3$, $-NR^8SO_2$ -phenyl, $-S(O)_2CF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$ -phenyl, or $-(CF_2)_rCF_3$;

each R^f is, independently at each occurrence, H, =O, OR^g , F, Cl, Br, I, CN, NO_2 , $-NR^8R^9$, $-C(O)R^g$, $-C(O)OR^g$, $-NR^8C(O)R^g$, $-C(O)NR^8R^9$, $-SO_2NR^8R^9$, $-NR^8SO_2NR^8R^9$, $-NR^8SO_2-C_{1-4}$ alkyl, $-NR^8SO_2CF_3$, $-NR^8SO_2$ -phenyl, $-S(O)_2CF_3$, $-S(O)_p-C_{1-4}$ alkyl, $-S(O)_p$ -phenyl, $-(CF_2)_rCF_3$, C_1-C_6 alkyl, C_2-C_6 alkenyl, or C_2-C_6 alkynyl;

each R^g is, independently at each occurrence, H, C_{1-6} alkyl, or $-(CH_2)_n$ -phenyl;

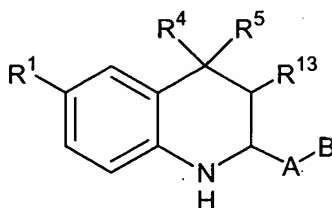
n, at each occurrence, is selected from 0, 1, 2, 3, and 4;

p, at each occurrence, is selected from 0, 1, and 2; and

r, at each occurrence, is selected from 0, 1, 2, 3, and 4;

provided that A is phenyl or a 6-membered aromatic N-heterocycle, then ring A is not substituted ortho to the tetrahydroquinoline with OH, halogen, $-CO_2H$, $-C(O)O-C_{1-4}$ alkyl, $-O$ -phenyl, $-O$ -benzyl, $-NR^7R^8$, $-CH_2OR^a$, haloalkyl, $-S-C_{1-4}$ alkyl, or $-NHSO_2-C_{1-4}$ alkyl.

3. (Currently amended) A compound according to Claim 2, wherein the compound is of Formula (Ib):



(Ib)

or a stereoisomer or a pharmaceutically acceptable salt or hydrate thereof, wherein:

B is phenyl substituted with 0-2 R^{11} and 0-1 R^{12} ;

R^1 is H, F, Cl, $-C(=NH)NH_2$, $-CH_2NH_2$, $-C(O)NR^7R^8$, OMe, or CN;

R^4 is H, $-(CH_2)_r-C_3-C_7$ cycloalkyl substituted with 0-2 R^{4b} , or $-(CH_2)_r$ -phenyl substituted with 0-3 R^{4b} , ~~or $-(CH_2)_r$ -5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^{4b} ;~~

each R^{4b} is, independently at each occurrence, H, OH, Cl, F, Br, CN, NO_2 , CF_3 , $-C(O)OR^a$, $-SO_2R^c$, $-NR^7R^8$, C_1 - C_4 alkyl, C_2 - C_4 alkenyl, C_2 - C_4 alkynyl, C_3 - C_6 cycloalkyl, C_1 - C_4 haloalkyl, C_1 - C_4 haloalkyloxy-, C_1 - C_4 alkyloxy-, C_1 - C_4 alkylthio-, C_1 - C_4 alkyl- $C(O)-$, or C_1 - C_4 alkyl- $C(O)NH-$;

R^5 is H, C_1 - C_3 alkyl, or C_3 - C_6 cycloalkyl;

each R^7 is, independently at each occurrence, H, C_{1-6} alkyl, or benzyl;

each R^{7a} is, independently at each occurrence, H, C_{1-4} alkyl substituted with 0-1 R^{7b} or 0-1 R^{7c} , $-(CH_2)_r$ - C_{3-7} cycloalkyl substituted with 0-1 R^f , or $-(CH_2)_r$ -phenyl substituted with 0-2 R^f , ~~or a $-(CH_2)_r$ -5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted 0-2 R^f ;~~

each R^{7b} is, independently at each occurrence, OR^g , F, CN, $-NR^7R^8$, $-C(O)R^g$, $-C(O)OR^g$, $-NR^8C(O)R^g$, $-C(O)NR^8R^9$, $-SO_2NR^8R^9$, or $-NR^8SO_2$ - C_{1-4} alkyl;

each R^{7c} is, ~~independently at each occurrence,~~ C_{3-7} cycloalkyl substituted with 0-1 R^f , or phenyl substituted with 0-2 R^f , ~~or a 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted 0-2 R^f ;~~

each R^8 is, independently at each occurrence, H, C_{1-6} alkyl, or benzyl;

each R^9 is, independently at each occurrence, H, C_{1-6} alkyl, or benzyl;

each R^{11} is, independently at each occurrence, H, F, $-(CH_2)_r$ - OR^a , CN, $-(CH_2)_r$ - NR^7R^8 , $-C(O)OR^a$, $-NR^8C(O)R^a$, $-NR^8C(O)OR^a$, $-C(O)NR^{7a}R^8$, $-NR^8C(O)NR^8R^9$, $-SO_2NR^8R^9$, or $-NR^8SO_2$ - C_{1-4} alkyl;

R^{12} is $-C(O)NR^{7a}R^8$, $-(CH_2)_r$ - CO_2R^{12a} , $-SO_2NHR^{12a}$, $-CONHSO_2NHR^{12a}$, $-SO_2NHCOR^{12a}$, $-SO_2NHCO_2R^{12a}$, $-CONHSO_2R^{12b}$, $-NHSO_2R^{12b}$, $-CONHSO_2R^{12b}$, $-CONHOR^{12b}$, or $-(CH_2)_r$ -5-tetrazolyl-;

each R^{12a} is, independently at each occurrence, H or C_{1-6} alkyl;

each R^{12b} is, independently at each occurrence, C_1 - C_4 alkyl substituted with 0-1 R^{12c} , C_2 - C_4 alkenyl substituted with 0-1 R^{12c} , C_2 - C_4 alkynyl substituted with 0-1 R^{12c} , - $(CH_2)_r$ - C_3 - C_7 carbocycle substituted with 0-2 R^{12c} , or - $(CH_2)_r$ -5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{12c} ;

each R^{12c} is, independently at each occurrence, H, F, Cl, Br, I, CF_3 , OCF_3 , CN, NO_2 , OR^a , $-CO_2R^a$, $-NR^7R^8$, $-SO_2R^c$, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, - $(CH_2)_r$ - C_3 - C_{10} carbocycle substituted with 0-3 R^d ; or - $(CH_2)_r$ -5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^d ;

R^{13} is H or C_1 - C_4 alkyl;

each R^a is, independently at each occurrence, H, C_{1-4} alkyl, $-(CH_2)_r$ - CO_2R^g , $-(CH_2)_r$ - C_{3-7} cycloalkyl, or $-(CH_2)_r$ - C_{6-10} aryl, ~~or $-(CH_2)_r$ -5-10 membered heteroaryl,~~
~~wherein said aryl or heteroaryl groups are substituted with 0-2 R^f ;~~

each R^f is, independently at each occurrence, H, =O, OR^g , F, Cl, Br, CF_3 , CN, NO_2 , $-NR^8R^9$, $-C(O)R^g$, $-C(O)OR^g$, $-NR^8C(O)R^g$, $-C(O)NR^8R^9$, $-SO_2NR^8R^9$, $-NR^8SO_2$ - C_{1-4} alkyl, $-NR^8SO_2CF_3$, $-S(O)_2CF_3$, $-S(O)_p$ - C_{1-4} alkyl, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, or C_2 - C_6 alkynyl;

each R^g is, independently at each occurrence, H or C_{1-4} alkyl;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, and 4; and

provided ring A is not substituted ortho to its attachment to the tetrahydroquinoline with OH, $-CO_2H$, $-C(O)O$ - C_{1-4} alkyl, O-phenyl, O-benzyl, $-NR^7R^8$, or $-NHSO_2C_{1-4}$ alkyl.

4. (Original) A compound according to Claim 3, wherein:

A is phenyl substituted with 0-2 R^{11} ;

B is phenyl substituted with 0-2 R¹¹ and 0-1 R¹²;

R¹ is -C(=NH)NH₂, -C(=O)NH₂, -CH₂NH₂, or OMe;

R⁴ is phenyl substituted with 0-1 R^{4b};

R^{4b} is H, OH, or F;

R⁵ is H, Me, Et, or Pr;

each R¹¹ is, independently at each occurrence, H, F, OH, OMe, CN, -NH₂, -CH₂OH, -CO₂H, -CO₂Me, -NHCOMe, -NHCOEt, -NHCOPr, -NHCO(i-Pr), -NHCO(i-Bu), -NHCO(cyclopropyl), -NHCO(phenyl), -NHCO(2-CO₂H-phenyl), -NHCO(3-CO₂H-phenyl), -NHCO(4-CO₂H-phenyl), -NHCO(3,5-(CO₂H)₂-phenyl)-, -NHCO(3,5-(CF₃)₂-phenyl), -NHCO(3-Me-5-CO₂H-phenyl), -NHCO(3-(t-Bu)-5-CO₂H-phenyl), -NHCO(3-CONH₂-5-CO₂H-phenyl), -NHCO(3-NH₂-5-CO₂H-phenyl), -NHCO(benzyl), -NHCO(phenethyl), -NHCO(phenylpropyl), -NHCO[2-(2-pyridyl)-ethyl], -NHCO(tetrazol-5-yl), -NHCOCH₂(tetrazol-5-yl), -NHCO(CH₂)₂(tetrazol-5-yl), -CONH₂, -CONHMe, -CONH(i-Pr), -CONH(i-Bu), -CONH(t-Bu), -CONH(benzyl), -CONH(phenethyl), -CONH(phenylpropyl), -CONH[2-(2-pyridyl)-ethyl], -NHCONHMe, -NHCONHEt, -NHCH₂CO₂H, -NHCOCO₂H, -NHCOCH₂CO₂H, -NHCO(CH₂)₂CO₂H, -NHCO(CH₂)₃CO₂H, -NHSO₂Me, -NHSO₂Et, or -CH₂NMe₂;

R¹² is -CO₂H, -CH₂(CO₂H), -CO₂Me, -SO₂NH₂, or -CONH₂;

R¹³ is H or Me; and

provided ring A is not substituted ortho to its attachment to the tetrahydroquinoline with OH, -CO₂H, -CO₂Me, -NH₂, or -NHSO₂C₁₋₄ alkyl.

5. (Original) A compound according to Claim 4, wherein:

A is 1,2-phenylene, 4-OMe-1,2-phenylene, 3-CO₂H-1,2-phenylene, 4-OMe-5-OH-1,2-phenylene, 5-CH₂OH-1,2-phenylene, 5-NHCOMe-1,2-phenylene, 5-phenylcarbamoyl-1,2-phenylene, 5-benzylcarbamoyl-1,2-phenylene, 5-phenethylcarbamoyl-1,2-phenylene, 5-(3-phenylpropylcarbamoyl)-1,2-phenylene,

5-[2-(2-pyridyl)ethylcarbamoyl]-1,2-phenylene, 5-NHCO(i-Bu)-1,2-phenylene, 1,3-phenylene, 6-OMe-1,3-phenylene, 6-F-1,3-phenylene, 5-NH₂-1,3-phenylene, 5-NHCOMe-1,3-phenylene, 5-NHCOEt-1,3-phenylene, 5-NHCOPr-1,3-phenylene, 5-NHCO(i-Pr)-1,3-phenylene, 5-NHCO(i-Bu)-1,3-phenylene, 5-NHCO(cyclopropyl)-1,3-phenylene, 5-NHCONHEt-1,3-phenylene, 5-NHCOCO₂H-1,3-phenylene, 5-NHCOCH₂CO₂H-1,3-phenylene, 5-NHCO(CH₂)₂CO₂H-1,3-phenylene, 5-NHCO(CH₂)₃CO₂H-1,3-phenylene, 5-NHCO(phenyl)-1,3-phenylene, 5-NHCO(benzyl)-1,3-phenylene, 5-NHCO(2-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(3-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(4-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(3,5-(CO₂H)₂-phenyl)-1,3-phenylene, 5-NHCO(3,5-(CF₃)₂-phenyl)-1,3-phenylene, 5-NHCO(3-Me-5-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(3-(t-Bu)-5-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(3-CONH₂-5-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(3-NH₂-5-CO₂H-phenyl)-1,3-phenylene, 5-NHCO(tetrazol-5-yl)-1,3-phenylene, 5-NHCOCH₂(tetrazol-5-yl)-1,3-phenylene, 5-NHCO(CH₂)₂(tetrazol-5-yl)-1,3-phenylene, 5-NHSO₂Et-1,3-phenylene, 5-NHCH₂CO₂H-1,3-phenylene, or 3-CO₂H-1,4-phenylene;

B is 2-CO₂H-phenyl, 4-CO₂H-phenyl, 2-SO₂NH₂-phenyl, 3-CH₂(CO₂H)-phenyl, 2,4-(CO₂H)₂-phenyl, 2,4-(CO₂Me)₂-phenyl, 2,4-(CONH₂)₂-phenyl, 2-CO₂H-4-CO₂Me-phenyl, 2-CO₂H-4-NH₂-phenyl, 2-CO₂H-4-CN-phenyl, 2-CO₂H-4-OMe-phenyl, 2-CO₂H-4-NHAc-phenyl, 2-CO₂H-4-CONH₂-phenyl, 2-CO₂H-4-CONH(i-Pr)-phenyl, 2-CO₂H-4-C(O)NH(i-Bu)-phenyl, 2-CO₂H-4-C(O)NH(t-Bu)-phenyl, 2-CO₂H-4-NHCOMe-phenyl, 2-CO₂H-4-NHCONHMe-phenyl, 2-CO₂H-4-CH₂NMe₂-phenyl, or 2-CO₂H-4-NHSO₂Me-phenyl;

R¹ is -C(=NH)NH₂, -C(=O)NH₂, -CH₂NH₂, or OMe;

R⁴ is phenyl, 4-OH-phenyl or 4-F-phenyl;

R⁵ is H, Me, Et, or Pr; and

R¹³ is H or Me.

6. (Original) A compound of Claim 1 selected from:

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-isobutylcarbamoyl-biphenyl-2-carboxylic acid;

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-isobutylcarbamoyl-5'-hydroxy-4'-methoxy-biphenyl-2-carboxylic acid;

2'-[6-carbamimidoyl-4-(4-hydroxy-phenyl)-1,2,3,4-tetrahydro-quinolin-2-yl]-5'-hydroxy-4-isobutylcarbamoyl-4'-methoxy-biphenyl-2-carboxylic acid;

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid dimethyl ester;

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2-carboxylic acid;

2'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-4-carboxylic acid;

2'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid dimethyl ester;

2'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-isobutylcarbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-t-butylcarbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2-carboxylic acid;

4-phenyl-2-(2'-sulfamoyl-biphenyl-3-yl)-1,2,3,4-tetrahydro-quinoline-6-carboxamide;

4-methyl-4-phenyl-2-(2'-sulfamoyl-biphenyl-3-yl)-1,2,3,4-tetrahydro-quinoline-6-carboxamide;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid 4-methyl ester;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid diamide;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

3'-(6-carbamimidoyl-4-ethyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

3'-(6-carbamimidoyl-4-propyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2,4-dicarboxylic acid;

4-amino-3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-(3-methyl-ureido)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-methanesulfonylamino-biphenyl-2-carboxylic acid;

4-acetylamino-3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-cyano-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-
biphenyl-2,4-dicarboxylic acid 4-methyl ester;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-
biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
methylcarbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
isopropylcarbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-t-
butylcarbamoyl-biphenyl-2-carboxylic acid;

3'-[6-carbamimidoyl-4-(4-fluoro-phenyl)-4-methyl-1,2,3,4-tetrahydro-quinolin-2-
yl]-4-carbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
dimethylaminomethyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-3-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
carbamoyl-biphenyl-2-carboxylic acid;

5'-amino-3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-
yl)-4-carbamoyl-biphenyl-2-carboxylic acid;

5'-amino-3'-(6-carbamimidoyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-
biphenyl-2-carboxylic acid;

5'-acetyl-amino-3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-
quinolin-2-yl)-4-carbamoyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
carbamoyl-5'-(3-methyl-butyrylamino)-biphenyl-2-carboxylic acid;

4-carbamoyl-3'-(6-carbamoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-
yl)-biphenyl-2-carboxylic acid;

4-carbamoyl-3'-(6-methoxy-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-
biphenyl-2-carboxylic acid;

3'-(6-aminomethyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-
carbamoyl-5'-(3-methyl-butyrylamino)-biphenyl-2-carboxylic acid;

3'-(6-aminomethyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-methyl-butyrylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-methylpropanoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(n-propanoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(cyclopropylcarbonylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-methoxyl-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(butyrylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-4'-methoxy-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-4'-fluoro-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-carboxyproacetyl amino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(carboxycarbonylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(benzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-methylpropanoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-phenylacetyl amino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-(4-fluorophenyl)-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-methylpropanoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-carboxypropanoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(4-carboxybenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-carboxybenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(2-carboxybenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(carboxymethylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3,5-bis(carboxybenzoylamino))-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-[(5-tetrazolyl)methylcarbonylamino]-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(4-carboxybutyrylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-[(5-tetrazolyl)carbonylamino]-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3,5-bisfluorobenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-amino-5-carboxybenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-[2-(5-tetrazolyl)ethylcarbonylamino]-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-carboxy-5-methylbenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-carboxy-5-t-butylbenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(3-aminocarbonyl-5-carboxybenzoylamino)-biphenyl-2-carboxylic acid;

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(ethylaminocarbonylamino)-biphenyl-2-carboxylic acid; and

3'-(6-carbamimidoyl-4-methyl-4-phenyl-1,2,3,4-tetrahydro-quinolin-2-yl)-4-carbamoyl-5'-(ethylsulfonylamino)-biphenyl-2-carboxylic acid;
or a stereoisomer or a pharmaceutically acceptable salt or hydrate thereof.

7. (Original) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 1 or a pharmaceutically acceptable salt or hydrate thereof.

8. (Original) A method for treating thromboembolic disorders, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1 or a pharmaceutically acceptable salt or hydrate thereof.

9. (Original) A method according to Claim 8; wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

10. (Original) A method according to Claim 9, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

11. (Withdrawn) A method for treating inflammatory disorders, comprising:
administering to a patient in need thereof a therapeutically effective amount of a
compound of Claim 1 or a pharmaceutically acceptable salt or hydrate thereof.
12. (Withdrawn) A method according to Claim 11, wherein the inflammatory disorder
is selected from the group consisting of sepsis, acute respiratory distress syndrome, and
systemic inflammatory response syndrome.
13. (Original) A method of treating a patient in need of thromboembolic disorder
treatment, comprising: administering a compound of Claim 1 or a pharmaceutically
acceptable salt or hydrate thereof in an amount effective to treat a thromboembolic
disorder.
14. (Original) A method, comprising: administering a compound of Claim 1 or a
pharmaceutically acceptable salt or hydrate thereof in an amount effective to treat a
thromboembolic disorder.
- 15-23. (Canceled)
24. (New) A pharmaceutical composition, comprising: a pharmaceutically
acceptable carrier and a therapeutically effective amount of a compound of Claim 2 or a
pharmaceutically acceptable salt or hydrate thereof.

25. (New) A method for treating thromboembolic disorders, comprising:
administering to a patient in need thereof a therapeutically effective amount of a
compound of Claim 2 or a pharmaceutically acceptable salt or hydrate thereof.

26. (New) A method according to Claim 25, wherein the thromboembolic disorder is
selected from the group consisting of arterial cardiovascular thromboembolic disorders,
venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the
chambers of the heart.

27. (New) A method according to Claim 26, wherein the thromboembolic disorder is
selected from unstable angina, an acute coronary syndrome, first myocardial infarction,
recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke,
atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein
thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral
arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and
thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling
catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other
procedures in which blood is exposed to an artificial surface that promotes thrombosis.

28. (New) A pharmaceutical composition, comprising: a pharmaceutically
acceptable carrier and a therapeutically effective amount of a compound of Claim 3 or a
pharmaceutically acceptable salt or hydrate thereof.

29. (New) A method for treating thromboembolic disorders, comprising:
administering to a patient in need thereof a therapeutically effective amount of a
compound of Claim 3 or a pharmaceutically acceptable salt or hydrate thereof.

30. (New) A method according to Claim 29, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

31. (New) A method according to Claim 30, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

32. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 4 or a pharmaceutically acceptable salt or hydrate thereof.

33. (New) A method for treating thromboembolic disorders, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 4 or a pharmaceutically acceptable salt or hydrate thereof.

34. (New) A method according to Claim 33, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders,

venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

35. (New) A method according to Claim 34, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

36. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 5 or a pharmaceutically acceptable salt or hydrate thereof.

37. (New) A method for treating thromboembolic disorders, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 5 or a pharmaceutically acceptable salt or hydrate thereof.

38. (New) A method according to Claim 37, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

39. (New) A method according to Claim 38, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

40. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 6 or a pharmaceutically acceptable salt or hydrate thereof.

41. (New) A method for treating thromboembolic disorders, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 6 or a pharmaceutically acceptable salt or hydrate thereof.

42. (New) A method according to Claim 41, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

43. (New) A method according to Claim 42, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein

thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.